

# Experimental Pauli-frame randomization on a superconducting qubit

Matthew Ware,<sup>1</sup> Guilhem Ribeill,<sup>1</sup> Diego Ristè,<sup>\*</sup> Colm A. Ryan,<sup>†</sup> Blake Johnson,<sup>‡</sup> and Marcus P. da Silva<sup>§</sup>

<sup>1</sup>*Raytheon BBN Technologies, 10 Moulton St., Cambridge, MA 02138, USA*

(Dated: March 23, 2021)

The promise of quantum computing with imperfect qubits relies on the ability of a quantum computing system to scale cheaply through error correction and fault-tolerance. While fault-tolerance requires relatively mild assumptions about the nature of qubit errors, the overhead associated with coherent and non-Markovian errors can be orders of magnitude larger than the overhead associated with purely stochastic Markovian errors. One proposal to address this challenge is to randomize the circuits of interest, shaping the errors to be stochastic Pauli errors but leaving the aggregate computation unaffected. The randomization technique can also suppress couplings to slow degrees of freedom associated with non-Markovian evolution. Here we demonstrate the implementation of *Pauli-frame randomization* in a superconducting circuit system, exploiting a flexible programming and control infrastructure to achieve this with low effort. We use high-accuracy gate-set tomography to characterize in detail the properties of the circuit error, with and without the randomization procedure, which allows us to make rigorous statements about Markovianity as well as the nature of the observed errors. We demonstrate that randomization suppresses signatures of non-Markovian evolution to statistically insignificant levels, from a Markovian model violation ranging from  $43\sigma$  to  $1987\sigma$ , down to violations between  $0.3\sigma$  and  $2.7\sigma$  under randomization. Moreover, we demonstrate that, under randomization, the experimental errors are well described by a Pauli error model, with model violations that are similarly insignificant (between  $0.8\sigma$  and  $2.7\sigma$ ). Importantly, all these improvements in the model accuracy were obtained without degradation to fidelity, and with some improvements to error rates as quantified by the diamond norm. This demonstrates the ability of Pauli-frame randomization to shape noise into forms that are more benign for quantum error correction and fault-tolerance.

## I. INTRODUCTION

Large-scale quantum computation poses a number of design and control challenges. Significant efforts are in progress [1–3] to meet and overcome challenges associated with initial state preparation, maintaining coherence, implementing universal gates, and measuring qubits reliably – all key criteria for building scalable quantum computers [4]. As the system coherence times continue to grow, coherent errors can become the dominant source of error. These errors can originate from miscalibration of qubit rotations, unintentional control frequency detunings, or interactions between systems that are otherwise assumed to be decoupled—all ubiquitous problems for experimental quantum computers. These errors are also particularly difficult to simulate in multiqubit systems, as they can interfere constructively and destructively, making prediction about the performance of quantum error correction codes and fault-tolerant computation quite difficult [5–7]. Moreover, theoretical lower bounds on the tolerable rates for coherent errors indicate they may be much more damaging than stochastic errors [8–11]. One way to address

this problem is to transform coherent errors into incoherent, stochastic errors, such as random bit and phase flips. Here we use a superconducting qubit system to implement *Pauli-Frame Randomization* (PFR) [12–15] and show that coherent errors can be reshaped into stochastic Pauli errors. We also discuss some additional benefits of the randomization process, such as decoupling of slow non-Markovian noise [16].

One significant challenge in determining whether PFR has indeed made coherent errors stochastic is their small magnitude. Thanks to the community’s progress towards fault-tolerance, the magnitudes of these errors are on the order of  $10^{-3}$  or less in state-of-the-art devices. Measuring such small errors reliably runs into limitations of various characterization approaches: standard tomography is sensitive to preparation and measurement imperfections and has very low accuracy, while randomized benchmarking estimates a quantity (closely associated with the average fidelity [17–19]) that does not differentiate between coherent and stochastic errors, and cannot test if errors corresponds to Pauli error models or not. In this demonstration we use *gate set tomography* (GST) [20–24], a tomographic reconstruction technique that provides 1) insensitivity to state preparation and measurement (SPAM) errors 2) nearly quantum-limited accuracy 3) an open source library for experiment design and data analysis [25]. Critically, GST also allows us to accurately quantify not only the behavior of the diamond norm error [26, 27] and average infidelity [17] under randomization, but also detailed features of individual gate errors and the degree to which the evolution is well described by a Markovian, time-invariant model [28]—all of which help confirm the

<sup>\*</sup> Current address: Quantum Engineering Solutions, Keysight Technologies, One Broadway, Cambridge, MA 02412

<sup>†</sup> Current address: AWS Center for Quantum Computing, Pasadena, CA 91125

<sup>‡</sup> Current address: IBM T.J. Watson Research Center, Yorktown Heights, NY 10598, USA; blake.johnson@ibm.com

<sup>§</sup> Current address: Microsoft Quantum, One Microsoft Way, Redmond, WA 98052; marcus.silva@microsoft.com

predicted Pauli error model behavior, despite the presence of general imperfections in the randomization operations.

The remainder of the paper is organized as follows. In Section II we describe PFR, and discuss how to test its implementation, in a statistically rigorous manner, in Section III. Section IV describes the experiments as well as the infrastructure required to create and process randomized sequences. Finally, in Section V we discuss the experimental results.

## II. PAULI-FRAME RANDOMIZATION

Pauli-frame randomization (PFR) is a noise-shaping technique that reduces general noise to effective random Pauli errors between computational gates [12–15]. If the computational gates consist of Clifford group operations [29] (a set of operations sufficient for the most promising approaches to error correction and fault-tolerance), the effect of these random Pauli operations can be easily tracked [30, 31] so that the computation can be unrandomized by simply reinterpreting the measurement results. While this randomization is designed to have no impact on the ideal computation, it effectively *symmetrizes* the error, much like *twirling* [32–37] and randomized decoupling [16], leading to an effective error operation that corresponds to a mixture of Pauli group operations known as a *Pauli channel*, or a *Pauli error model*.

These results can be derived in the limit of perfect randomization operations and gate-independent errors as follows. Consider a set of ideal (resp. noisy) [38] Clifford group quantum operations  $C_i$  (resp.  $\tilde{C}_i$ ). Any sequence of ideal Clifford operations can be randomized by inserting uniformly random Pauli group operations between the Clifford group operations. Since Clifford group operations transform Pauli group operation to other Pauli group operations, the overall effect of these random Pauli group operations can be cancelled out by applying a final single Pauli group operation at the end of the sequence of gates. Moreover, since the Pauli group is a subgroup of the Clifford group, one may simply combine the  $i$ th random Pauli operation  $P_i$  with the  $i$ th Clifford group operation  $C_i$ , to obtain a random Clifford group operations  $D_i$  [39]. In other words, a given sequence of Clifford group operations  $C_L C_{L-1} \cdots C_2 C_1$  becomes

$$\underbrace{P_{L+1} C_L P_L}_{\mathcal{D}_L} \underbrace{C_{L-1} P_{L-1}}_{\mathcal{D}_{L-1}} \cdots \underbrace{C_2 P_2}_{\mathcal{D}_2} \underbrace{C_1 P_1}_{\mathcal{D}_1} \quad (1)$$

which results in the randomized sequence of Clifford group operations  $\mathcal{D}_L \mathcal{D}_{L-1} \cdots \mathcal{D}_2 \mathcal{D}_1$ . In essence, under PFR, a single realization of a randomized sequence of Clifford group operations simply corresponds to a different sequence of Clifford group operations.

It is possible to choose all  $P_i$  independently at random and compensate for their action by flipping observed measurement outcomes in post-processing (as, by construction, we only measure in the computational basis).

In order to simplify post-processing, we instead choose  $P_{L+1}$  to cancel the effect that all other random Pauli group operations would have on measurement results (i.e.,  $P_{L+1}$  is a Pauli frame correction before measurement). In this way the measurement outcome of the randomized and unrandomized experiments can be treated exactly the same, with no additional post-processing for the randomized experiments.

We can analyse the sequences above with the simplifying assumption of gate-independent errors by replacing each operation with its noisy counterpart. We write the noisy operations  $\tilde{D}_i = \mathcal{E} D_i$  (where  $\mathcal{E}$  is an arbitrary but fixed completely-positive trace-preserving (CPTP) map) to obtain

$$\tilde{D}_L \tilde{D}_{L-1} \cdots \tilde{D}_2 \tilde{D}_1, \quad (2)$$

$$= \mathcal{E} D_L \mathcal{E} D_{L-1} \cdots \mathcal{E} D_2 \mathcal{E} D_1, \quad (3)$$

$$= \mathcal{E} P_{L+1} C_L P_L \mathcal{E} C_{L-1} P_{L-1} \cdots \mathcal{E} C_2 P_2 \mathcal{E} C_1 P_1. \quad (4)$$

Defining  $\mathcal{P}^C = C P C^\dagger$ , we can write  $C P = \mathcal{P}^C C$ . Similarly, we define  $\mathcal{P}_{n:1} = P_n \mathcal{P}_{n-1:1}^{C_{n-1}}$  (with the base case  $\mathcal{P}_{1:1} = P_1$ ). With these definitions, the entire sequence can then be rewritten as  $\mathcal{E} P_{L+1:1} C_L \mathcal{P}_{L-1:1}^{C_{L-1}} \mathcal{E} \mathcal{P}_{L-1:1}^{C_{L-1}} C_{L-1} \cdots \mathcal{P}_{1:1}^{C_1} \mathcal{E} \mathcal{P}_{1:1}^{C_1} C_1$ , where, in the experiments described here, we have chosen  $P_{L+1:1}$  to be the identity. In other words, we choose  $P_i$  uniformly at random for  $1 \leq i \leq L$ , and choose  $P_{L+1}$  to get a trivial  $\mathcal{P}_{L+1:1}$ . Averaging over many uniformly random choices of Pauli operations in Eq. 4, we transform each  $\mathcal{E}$  in the sequence into  $\bar{\mathcal{E}} = \frac{1}{d^2} \sum_i P_i \mathcal{E} P_i$ , which correspond to twirling  $\mathcal{E}$  over the Pauli group. This, in turn, ensures that the effective error  $\bar{\mathcal{E}}$  associated with each gate in the sequence corresponds to a statistical mixture of Pauli operations [35], as desired [40].

The calculation outlined above does require rather strong assumptions about the properties of the noise (i.e., that it is gate independent and Markovian), but due to similarities to randomized benchmarking (RB) [41–45], which has been shown to require weaker assumptions [19], we expect that these strong assumptions are not strictly necessary. In the remainder of this paper we focus on how to test such a hypothesis, and implement these test on the natural imperfections of a superconducting qubit experiment.

## III. HYPOTHESIS TESTING

The task of checking whether the result of applying PFR to an experiment does indeed result in a Pauli channel is subtle. Modern experiments have very high fidelity to ideal operations so checking that the unrandomized errors are not well described by Pauli channel—i.e., determining that PFR is necessary—is already challenging, since error rates can be on the order of  $10^{-3}$  or less. In both cases, it is natural to consider long sequences of operations to amplify sensitivity to these small errors.

We choose to use long-sequence gate-set tomography (GST) [22–24] to observe these small effects, and use a readily available open-source package for experiment design and data analysis [25], with minor modifications. At heart, GST is a sophisticated refinement of a quantum process tomography [46, 47], providing a complete reconstruction of the action of quantum operations. In particular, GST is an iterative procedure that refines the tomographic reconstruction of a set of gates by comparing predictions about long gate sequences to experimental observations, and adjusting the reconstruction for better agreement. Since long sequences allow for small perturbations to accumulate, this technique yields unparalleled accuracy [22–24, 48].

Even with a reconstruction in hand, another subtle question is how to quantify the distance between reconstructed errors and a Pauli error model—i.e., the degree of “non-Pauliness” of the noise. We use the likelihood ratio test for this purpose [49, 50], which requires a hierarchy of nested models. The null hypothesis  $H_0$  is taken to be that the statistics for each sequence in the GST experiments leads to a separate Binomial distribution of outcomes. More explicitly, for the null hypothesis we only assume that the sequences correspond to reproducible experiments with well defined measurement statistics, and ignore the gate structure of the sequences. This corresponds to not making any assumption about Markovianity or time independence of the system evolution. We then consider two hypotheses nested within  $H_0$ : that each gate in the sequence corresponds to a fixed linear operation acting on the system (we call this first hypothesis  $H_1$ ), and that each gate in the sequence corresponds to a fixed Clifford group operation followed by a fixed Pauli stochastic error operation (we call this second hypothesis  $H_2$ ). The consistency of  $H_0$  with  $H_2$ , for some reconstructed Pauli error model, will be taken as our measure of Pauliness.

As we indicated, these hypotheses are nested:  $H_2$  is a special case of  $H_1$ , and  $H_1$  is a special case of  $H_0$ , meaning that if the statistical tests indicated  $H_2$  is consistent with  $H_0$ , the same will be true of  $H_1$ . The statistical tests will only be able to test if the proposed hypotheses are consistent with  $H_0$ , in the sense that a hypothesis cannot have a higher likelihood than another hypothesis it is nested into.

We fit data to a model under  $H_0$  by maximum-likelihood estimation of the Binomial distribution parameter  $p$  associated with each GST sequence. We fit data to a model under  $H_1$  using progressive refinement of maximum-likelihood estimation, a heuristic developed for GST [25]. We fit data to a model under  $H_2$  by projecting the fit of  $H_1$  into a generalized monomial matrix (described below), determined by the corresponding noiseless Clifford group operation. The first two fits are part of the standard routines within GST, while the last fit is a small extension to the existing GST routines.

The fitting of data to a model under  $H_2$  proceeds as follows. In the Pauli-Liouville representation [51–57], a Clifford group operation is a monomial matrix—each row

or column has a single non-zero matrix element, and this matrix element is  $\pm 1$ . In the presence of a Pauli error model, a noisy Clifford group operation will be a generalized monomial matrix, where the  $\pm 1$  elements of the noiseless matrix are replaced by numbers in the interval  $[-1, 1]$  (but the 0 matrix elements remain unchanged). Collectively, these matrix elements must live in a simplex equivalent to the probability simplex for the Pauli channel [36]. Thus, the projection of an  $H_1$  model onto an  $H_2$  model simply corresponds to identifying which matrix elements should be set to zero (i.e., which matrix elements are zero in the ideal gate), and then adjusting the remaining non-zero matrix elements so that the resulting matrix lies in the appropriate simplex.

### Badness-of-fit

We quantify how well the data is explained with each of the hypotheses discussed above by computing a metric for the quality of the fits obtained. The basis for this calculation is  $\mathcal{L}(H_i)$ , the likelihood of the observed data given the model fitted under a particular hypothesis  $H_i$ .

Following Wilk’s theorem [50], we know the log-likelihood ratio  $-2 \log \frac{\mathcal{L}(H_i)}{\mathcal{L}(H_0)}$  has a distribution that asymptotically (in the same size) approaches a  $\chi^2$  distribution with degrees of freedom given by the difference in the dimensionality of the two nested hypotheses, under the assumption that the null hypothesis is true. The mean and variance of the asymptotic distribution for the log-likelihood ratio are determined by the number of degrees of freedom. Given the fitted models, the likelihood of the observations under the various hypotheses are computed, and we follow the convention of reporting the difference between the observed statistic and the mean predicted by Wilk’s theorem, in units of the standard deviation of the appropriate  $\chi^2$  distribution, and call this quantity  $N_\sigma$ . Intuitively, if this “badness-of-fit” number is large, we favor the null hypothesis (i.e., the hypothesis fit is bad), but if this number is small, both the simpler hypothesis and the null hypotheses are valid, and the simpler hypothesis is favored as a parsimonious model. We emphasize this “badness-of-fit” parameter cannot be obtained by characterization techniques like RB that do not have an implicit model built in.

The log-likelihood ratios allow us to quantify whether (a) the observations are consistent with a Markovian error model (i.e., whether  $H_1$  is plausible), and (b) whether the observations are consistent with a Clifford group operation with a Pauli error model (i.e., whether  $H_2$  is plausible). In particular, we are interested in testing whether the answer to these questions changes when we apply PFR to our experiments. For this, it is necessary to look at the likelihood of hypotheses in different data sets (i.e., the unrandomized and the randomized GST experiments). Likelihoods cannot be meaningfully compared across different data sets. Instead, we simply consider the plausibility of the different hypothesis for the differ-

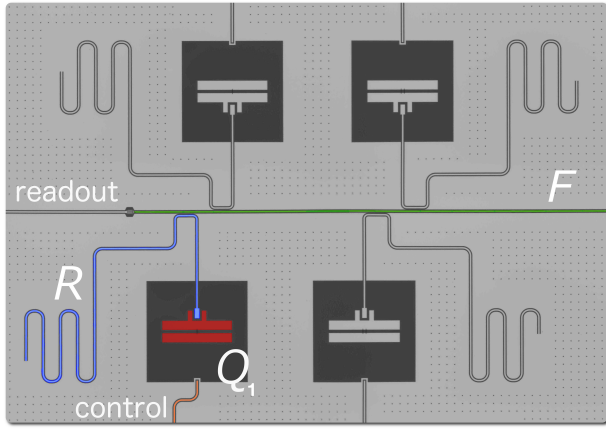


FIG. 1. False-color micrograph showing qubit  $Q_1$  (red), resonator  $R$  (blue) and Purcell filter  $F$  (green). The qubit is dispersively coupled to a  $\lambda/4$  readout resonator which is capacitively coupled to a Purcell filter with a  $Q = 22$ . Qubit control is done through a dedicated drive line (orange) and all qubit readout is done via a central feed line incorporating a Purcell filter.

ent data sets, while taking great care to ensure that the data sets are representative of the same noise and error environment, as we now describe.

## IV. EXPERIMENTS

### Device parameters

To test the hypotheses of Section III, we implement the PFR procedure on a superconducting qubit device. The device consists of four fixed-frequency, transmon qubits, designed to be similar to those described in [3]. The qubits are uncoupled but readout through a common Purcell filter. For the PFR experiment in this Letter, only one qubit ( $Q_1$ ) is measured.  $Q_1$  is dispersively coupled to a readout resonator with a center frequency of  $\omega_r/2\pi = 7.112$  GHz,  $\kappa/2\pi = 3.4$  MHz, which is in turn capacitively coupled to a quarter-wave Purcell filter with external  $Q = 22$  and a center frequency of  $\omega_f = 7.27$  GHz [58] enabling fast qubit readout.  $Q_1$  has a fixed 0-1 transition frequency of  $\omega_q/2\pi = 4.432$  GHz with an anharmonicity  $\alpha/2\pi = 308$  MHz. Coherence times measured for  $Q_1$  are  $T_1 = 10 \mu\text{s}$ ,  $T_2 = 13 \mu\text{s}$  and a Hahn echo time of  $T_{\text{echo}} = 16 \mu\text{s}$  (the other qubits in the device were not characterized). The pulses used were 50ns long, leading to an expected average gate infidelity (resp. diamond norm distance) of at least  $\sim 0.2\%$  per pulse ( $\sim 1\%$  per pulse). Since we use 2 pulses in the implementation of the gates discussed here, we expect an infidelity of no less than  $\sim 0.4\%$  and diamond norm distance of no less than  $\sim 2\%$  (ignoring all other sources of error and ignoring the effects of PFR).

### Electronics and software stack

Making the PFR process experimentally tractable requires leveraging a complex software and hardware control infrastructure. The first hurdle is the sheer number of experiments needed. Long-sequence GST ( $\ell\text{GST}$ ) experiments require a large set ( $\sim 3500$ ) of long circuits, each with up to 6155 gates. To ensure high accuracy, we produce a large number of measurement shots, 1000 per sequence. Under PFR, we take a single measurement shot per randomized sequence, resulting in 1000 unique GST circuits for each of the  $\sim 3500$   $\ell\text{GST}$  circuits originally specified. Thus, in total, we measure over 3.5 million unique sequences to obtain high tomographic reconstructions for the gates in the unrandomized and the randomized experiments [59]. The second challenge is running the experiments in a way that allows the most direct comparison between the randomized and unrandomized cases—doing so allows us to minimize the impact of drift when comparing how the hypothesis tests from Sec. III fare on the different data sets. To achieve this, the unrandomized and the randomized sequences should be run in an interleaved fashion to ensure they experience the same noise environment (to the extent possible). These requirements necessitate hardware that can execute a large number of very long circuits, and to quickly alternate between them.

To address these issues we use a custom sequence compiler written in Julia [60] called `QGL.jl` [61], providing a  $4\times$  compilation speed-up per circuit over an earlier Python version through a combination of parallel computation and other efficiency improvements—this ensured we were able to compile the 3.5 million unique sequences into pulse sequences in a reasonable amount of time. To minimize the runtime overhead we leverage a custom arbitrary pulse sequencer with gateway dedicated to implementing quantum circuits [62].

A rough outline of the process is as follows. Standard, one-qubit  $\ell\text{GST}$  sequences are created using the GST experiment and analysis software `pyGSTi` [25]. For the data presented here, we choose the maximal sequence length in GST to be 6150 gates, to ensure the experiment will have high accuracy, and be sensitive to non-Markovianity over timescales long compared to qubit coherence times. The  $\ell\text{GST}$  sequences are then randomized as described in Sec. II. It is worth emphasizing the lengths of the randomized circuits are unchanged as the Pauli group operations are combined with a neighboring Clifford group operation [63], much like the randomized compiling proposal of Ref. [15].

The collection of uniquely randomized GST (rGST) circuits and the original unrandomized circuits are then passed to the `QGL.jl` compiler which translates qubit gate instructions into machine instructions. This involves not only mapping high-level instructions to control pulses but also time-ordering and synchronizing instruction playback between all qubit control and readout channels. We note, this process takes significantly more time than the genera-



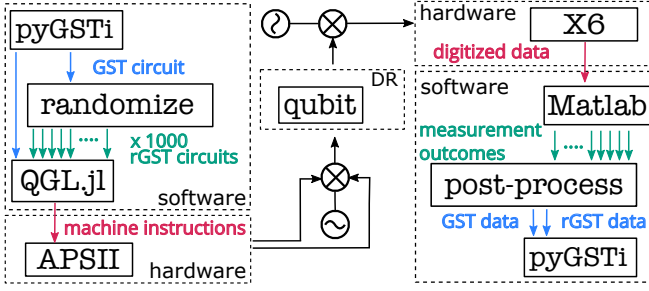


FIG. 2. Experimental data flow. Basic GST sequences are created using pyGSTi [25]. This basic experiment is then randomized 1000 times as described in section II. Each randomization and the original experiment get compiled by the QGL.jl compiler that translates sequence instruction into instructions implemented by APSII pulse sequencer [64]. The qubit response is digitized by an Innovative Integration X6 digitizer card and organized with a Matlab experimental framework. The single-shot data from each experiment is then post-processed into counts that pyGSTi uses to reconstruct the gate set process matrices and the goodness-of-fit metrics.

tion or randomization of the GST circuits. The compiled instructions are then passed to the pulse sequencer which is used to control the qubit. Due to the nature of the randomization process, the rGST experiments lack any kind of repetition or subroutine structure which rules out any efficient storage in hardware memory of a complete set of circuits. To address this, the set of randomized experiments are broken into groups of 10 in order to fit in the control hardware memory. These 10 single shot runs of rGST were then interleaved with 10 shots of unrandomized  $\ell$ GST. The process is repeated 100 times for each data point in Fig. 3. The complete process flow for a single round of experiment generation is illustrated in Fig. 2.

In the canonical construction of the Clifford group, elements are composed of multiple native  $\pi$  and  $\pi/2$  pulses, which leads to Clifford group elements being implemented by different numbers of native gates/pulses non-uniform length. To account for this, we use a “diatomic” implementation of the group where each Clifford group operation is performed with two  $X_{\pi/2}$  pulses of fixed length (50 ns) and three possible Z-frame updates [57, 65]. This diatomic approach ensures all Clifford operations have equal duration. The room temperature measurement signals were processed with an autodyne technique described in Ref. [66] using the BBN-QDSP digitization architecture [64] for the *Innovative Integrations X6-1000M* digitizer card. The final state assignment is then fed into the pyGSTi package for gate set reconstruction. pyGSTi also provides the likelihood of  $H_0$  and  $H_1$ , while custom code generates the likelihood of  $H_2$ —from these likelihoods, we obtain the likelihood ratio statistic and compare it to the predictions from Wilk’s theorem.

## V. RESULTS

The experiment outlined in Sec. III was performed to test the effectiveness of PFR. This process was repeated seven times, each taking roughly one hour to complete. The repetitions allow us to observe how drift affects the results over an operationally meaningful amount of time.

One of the critical questions of this work is the validity of  $H_2$  (the hypothesis that gates are well described by Clifford group operations followed by stochastic Pauli noise) and the Markovian behavior of qubit evolution under PFR. Data addressing this question can be seen in Fig. 3 where the GST model violation is plotted in terms of  $N_\sigma$  both with and without randomization. Several features are immediately apparent: (1) the Markovian fits ( $H_1$ ) to the unrandomized experiments (filled triangles) are orders of magnitude worse than the randomized experiments (empty triangles), (2) the data projected to a Pauli error model ( $H_2$ ) in the unrandomized cases (red filled triangles pointing down) is roughly three orders of magnitude worse than the randomized experiments (red empty triangles pointing down), (3) there is little difference between the quality of the fits under all the hypotheses for the randomized experiments (empty triangles). In terms of the hypotheses outlined previously, for unrandomized experiments there is a large likelihood discrepancy between  $H_0$  and the simpler hypotheses, greatly favoring the non-Pauli, non-Markovian  $H_0$  model ( $H_1$  is  $43\sigma$  to  $76\sigma$  away from the predictions from Wilk’s theorem, and  $H_2$  is  $1754\sigma$  to  $1987\sigma$  away), while for the randomized experiments all hypotheses have comparable likelihood (within  $0.3\sigma$  to  $2.7\sigma$  of the predictions from Wilk’s theorem), so it is reasonable to take the simplest hypothesis (the Markovian, stochastic Pauli error model  $H_2$ ) as the best explanation for those observations.

We should note that, despite the base level of  $H_1$  model violation measured in the unrandomized data (a signature of non-Markovianity) appearing large at  $1987\sigma$ , it is largely consistent with observations in other systems under similar circumstances (see, e.g., ion trap experiments without drift control or decoupling pulses [23]).

These features strongly indicate the noise in the absence of randomization is not well described by a Markovian error model, which follows from comment (1) above. Also apparent from comment (2) is that even the best Markovian error model is not well approximated by a Pauli model in the absence of randomization. Conversely, these features indicate the noise under PFR is very well described by a Markovian Pauli error model. In much simpler terms, the features of non-Pauli error models (i.e., non-trivial off-diagonal matrix elements in the Pauli-Liouville representation [67]) are insignificant in the reconstructions of the randomized experiments, as Fig. 4 illustrates. These separations are persistent over many repeats of the experiment, and the separation of many orders of magnitude indicates that PFR worked in these experiments not only quantitatively, but qualitatively, i.e., unrandomized experiments have strong non-Markovian fea-

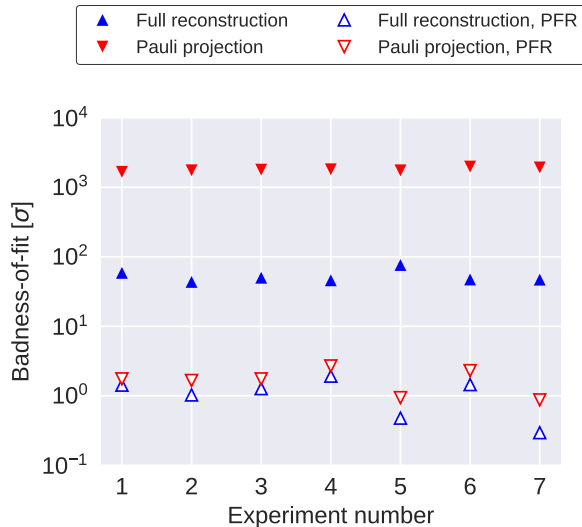


FIG. 3. Badness-of-fit for the GST reconstructions under a Markovian error model ( $H_1$ , upward blue triangles), and a Markovian stochastic Pauli error model ( $H_2$ , downward red triangles), as quantified by the log-likelihood-ratio statistic. This statistic is presented as the difference from the predicted mean of the  $\chi^2$  distribution (from Wilk’s theorem), in units of the standard deviation of that same distribution, under the assumption that  $H_0$  is true. Both experiments without randomization (full triangles) and with randomization (empty triangles) are considered. As discussed in the text, the 3.5 million unique sequences that comprise the tomography experiments for randomized and unrandomized experiments are measured in an interleaved fashion, so that both reconstructions should experience the same physical noise conditions. The entire collection of tomography experiments are repeated 7 times to illustrate the behavior observed persists, and thus unlikely to be the result of statistical fluctuations, and is robust to drift in our system (each of these 7 experiment lasted roughly one hour).

tures, while randomized experiments were well explained by Markovian Pauli error models.

### Behavior of error metrics under randomization

The badness-of-fit results illustrate that the models derived under randomization are much more useful in explaining the observations than the models derived without randomization. A natural question that arises is whether this comes at the cost of degrading the performance of the gates. Here we demonstrate that this is not the case, and that in fact the performance of the gates improves under PFR.

We computed the average gate infidelity [69] and the diamond norm distance [70] for the reconstructed gates under normal operation and under PFR, as depicted in Fig. 5. The observed average gate infidelities (resp. di-

among norm distances) are roughly double (quadruple) the expected coherence limits of the device, which may be explained by dynamical effects in the gate implementation which may be addressed by more careful pulse shaping [71] (and which are not accounted for in the coherence limit calculation mentioned earlier). We observe no appreciable difference between the infidelity of randomized and unrandomized experiments, while the diamond norm distance is reduced by a factor of 3-5 under PFR. This is consistent with the well known behavior of the infidelity and the diamond norm under small coherent errors—namely, the infidelity is only sensitive to coherent errors to second order, while the diamond distance is sensitive to first order [72].

The diamond norm distance of unrandomized experiments monotonically increases over the course of the 7 experiments, a behavior consistent with drift in the qubit and control parameters with respect to calibrations. The qubit control parameters are calibrated only once, at the beginning of the first experiment. This drift may at least partially explain the violation of the time-independent Markovian model represented by  $H_1$ , since these parameters appear to be continuously and systematically drifting throughout the 7 experimental runs, but this is a small effect since  $H_1$  still makes accurate predictions within each of the runs.

It should be noted that the drift is not apparent in the randomized experiments (even in the diamond norm distance), despite these experiments being run under the same conditions as the unrandomized experiments. This indicates that the drift was averaged away under PFR, a behavior consistent with coherent errors.

### Access to data

The experimental data, along with scripts used to perform the analysis and plot the results, can be found in Ref. [68]. These include the full tomographic reconstruction of gatesets for all 7 experiments, along with the raw counts needed for these reconstructions.

## VI. SUMMARY

We have demonstrated that Pauli-frame randomization reduces both the non-Markovian features and the non-Pauli model features of errors in single qubit experiments. This demonstration relies on long-sequence gate-set tomography, which yields high accuracy reconstructions of all operations used in the experiments. This in turn required a high-degree of automation to capture and process the  $\sim 7$  million measurement shots/hour. In the absence of randomization, the experiments were shown to have strong non-Markovian features, and the best Markovian model in that case was also shown to have strong features inconsistent with Pauli error models. In the presence of Pauli-frame randomization, the experiments were shown

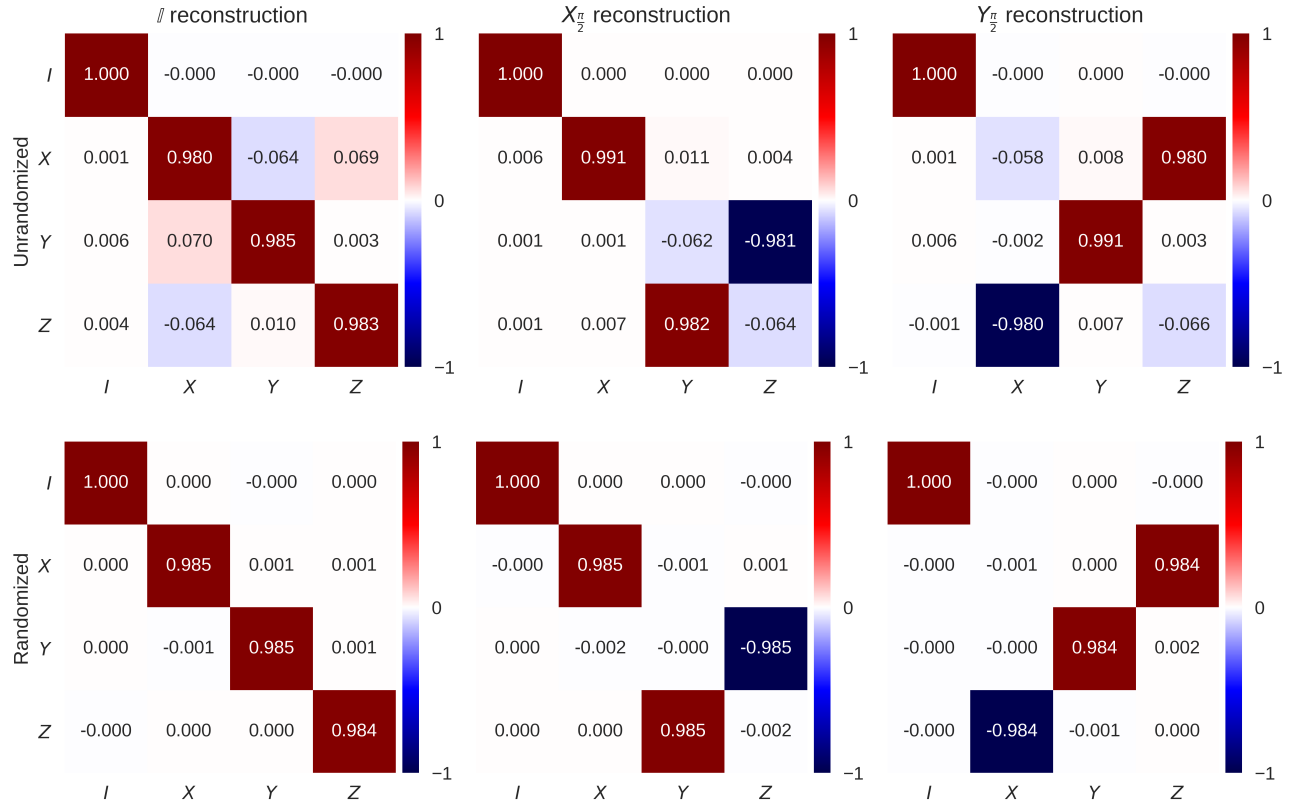


FIG. 4. Matrix representations of the reconstructed processes corresponding to the  $I$ ,  $X_{\pi/2}$ , and  $Y_{\pi/2}$  operations in the first of the seven experiments performed (details of the other six experiments and confidence interval computation can be found in Ref. [68]). Without randomization (top row) there are significant off-diagonal contributions, corresponding to non-Pauli errors. With randomization (bottom row) there are no statistically significant off-diagonal contributions, indicating the errors correspond to a Pauli error model, as expected. Error bars (95% confidence) for this experiment are smaller than  $\pm 0.0028$  (unrandomized) and  $\pm 0.0023$  (randomized).

to be highly consistent with a Markovian Pauli error model, as predicted. As quantified by log-likelihood-ratio statistic, the violation of Markovian and Pauli error models in the unrandomized experiments is highly-significant, as high as  $1987\sigma$ , while the violation of Markovian Pauli error models in the randomized experiments are statistically insignificant, less than  $2.7\sigma$  in most of the experiments. This several orders-of-magnitude separation between randomized and unrandomized experiments was persistent across seven repeats of the experiment, indicating the noise-shaping effect of Pauli-frame randomization is robust to drift in the control parameters and fluctuations in the noise environment.

Areas for future work include speeding up the experiments using techniques such as active reset [73, 74], and pushing randomization process onto the hardware FPGA, which would allow for data acquisition of randomized Clifford group circuits without the user having to manually pre-compile random circuits.

After this work was completed similar results were independently reported by Hashim et al. [75].

## VII. ACKNOWLEDGEMENTS

We thank Robin Blume-Kohout, Erik Nielsen, Joel Wallman, and Joseph Emerson for fruitful discussions, and the anonymous referees for comments that helped improved the presentation of the paper. We also thank Alan Howsare, Adam Moldawer and Ram Chelakara at Raytheon Integrated Defense Systems for sample fabrication.

The qubit design and fabrication were funded by Internal Research and Development at Raytheon BBN Technologies and directed by Thomas Ohki. This work was funded by LPS/ARO grant W911NF-14-C-0048. The content of this paper does not necessarily reflect the position or the policy of the Government, and no official endorsement should be inferred.

## CONTRIBUTIONS

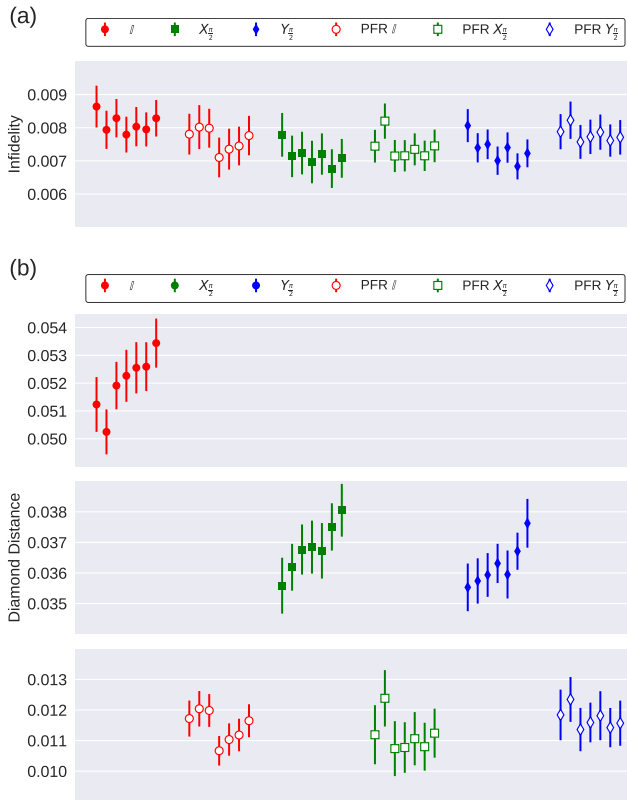


FIG. 5. (a) Average gate infidelity and (b) diamond norm distance estimates for all three gates (colors) on a sequence of seven separate experiments over several hours. Data for different gates is horizontally offset for clarity, but we emphasize each of the experiments leads to a reconstruction of all three gates, for both randomized (full symbols) and unrandomized (empty symbols) gate-sets under identical conditions—the only distinction between the seven experiments is the passage of time. All quantities are computed from the reconstructed process matrices. For the randomized case (empty symbols), the infidelity and the diamond norm are comparable, at  $\approx 1\%$ . For the unrandomized experiments, there is significant deviation between the diamond norm error rate and the infidelity, suggesting the presence of coherent errors that affect the infidelity metric only weakly (and which are suppressed in the randomized experiments). A monotonic upward trend in the diamond norm distance of unrandomized experiments (full symbols) implies the presence of systematic drift in the control pulses, which is also suppressed by randomization (empty symbols). Error bars are 95% confidence intervals (analytically for (a), and with the Hessian provided by pyGSTi for (b)).

MW, GR and MPS wrote the manuscript. MW, GR, and DR performed the experiment, while CR and BJ built the control software and electronics infrastructure. GR and MW designed the device and assisted in fabrication. MPS proposed the experiment, and wrote the experiment generation scripts. MW and MPS performed the data analysis.

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$$\mathcal{F}(\mathcal{U}, \tilde{\mathcal{U}}) = \frac{\text{tr } \tilde{\mathcal{U}}\mathcal{U}^\dagger + d}{d^2 + d}$$

- [70] The diamond distance between two processes [26, 27] is defined as

$$\mathcal{D}_\diamond = \frac{1}{2} \|\tilde{\mathcal{U}} - \mathcal{U}\|_\diamond = \sup_\rho \frac{1}{2} \|(\tilde{\mathcal{U}} \otimes \mathcal{I}_d - \mathcal{U} \otimes \mathcal{I}_d)(\rho)\|_1,$$

with  $d$  being the total system dimension.

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